

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080919\
 Data File : BF116124.D
 Acq On : 9 Aug 2019 20:39
 Operator : HP/JU
 Sample : K4244-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BERGEN-SP-1

Manual Integrations
 APPROVED

Sohil
 8/13/2019 7:18:01 AM

Quant Time: Aug 10 05:35:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 11:13:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	131472	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	511554	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	265138	20.00	ng	-0.02
64) Phenanthrene-d10	11.36	188	462315	20.00	ng	-0.01
76) Chrysene-d12	14.00	240	305541	20.00	ng	-0.01
87) Perylene-d12	15.46	264	336483	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	784370	99.88	ng	0.00
7) Phenol-d6	6.47	99	999517	100.87	ng	-0.01
23) Nitrobenzene-d5	7.40	82	639918	72.21	ng	-0.02
42) 2,4,6-Tribromophenol	10.66	330	253381	98.57	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	1085922	76.72	ng	-0.01
79) Terphenyl-d14	12.94	244	1086033	74.90	ng	-0.01
Target Compounds						
10) Phenol	6.49	94	40714	3.489	ng	92
50) Dimethylphthalate	9.59	163	102381	5.671	ng	98
75) Fluoranthene	12.57	202	51268	2.072	ng	98
78) Pyrene	12.80	202	57560	2.499	ng	98
88) Benzo(b)fluoranthene	15.04	252	52188m	2.682	ng	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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